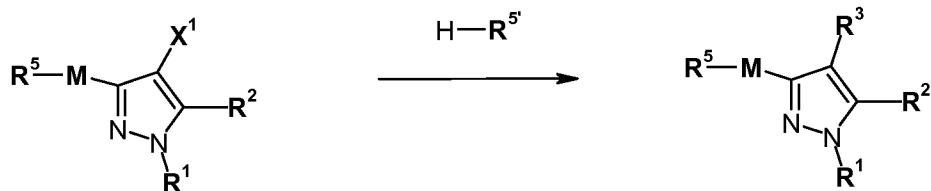


Listings of claims:

- 1-10. (Canceled)
11. (Currently Amended) A compound 2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-(1,3-benzodioxol-5-yl)ethyl]-(2*S*)-propylamine or a salt, or in-vivo hydrolyzable ester ~~pro-drug or solvate~~ thereof.
12. (Cancelled)
13. (Currently amended) A pharmaceutical formulation comprising a compound, or salt, or in-vivo hydrolyzable ester ~~pro-drug or solvate~~ thereof, according to claim 11 and a pharmaceutically acceptable diluent or carrier.
14. (Withdrawn) A method of antagonising gonadotropin releasing hormone activity in a patient, the method comprising administering a compound, or salt, pro-drug or solvate thereof, according to claim 1 to a patient.
15. (Withdrawn) A method of treating and/or preventing a sex hormone related condition in a patient, the method comprising administering a compound according to claim 1, or salt, pro-drug or solvate thereof, to a patient.
16. (Withdrawn) A process for the preparation of a compound of Formula (I) as defined in Claim 1, comprising a process selected from (a) to (h) as follows:
- (a) Reaction of a compound of formula **XXXII** with a compound of formula **H-R⁵**, to form a compound of Formula (I),



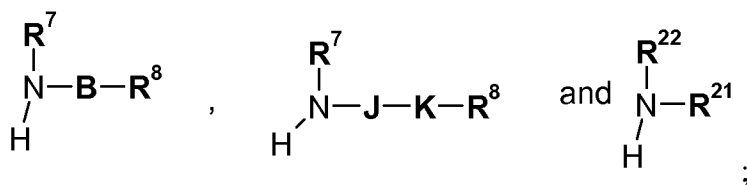
XXXII

Formula (I)

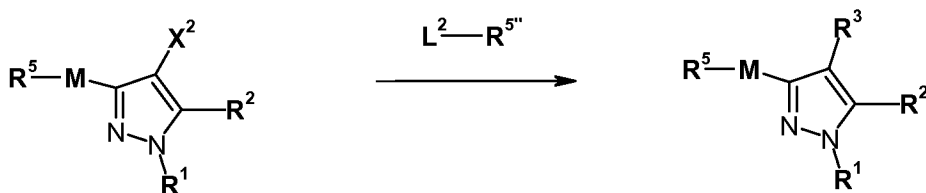
wherein X^1 is selected from:

and L^1 is a displaceable group; and

$H-R^{5'}$ is selected from:



- (b) Reaction of a compound of formula XXXIII with a compound of formula $H-R^{5''}$ to form a compound of Formula (I),



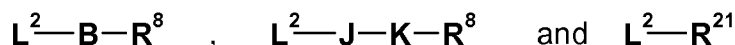
XXXIII

Formula (I)

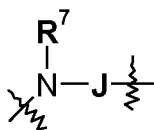
wherein X^2 is selected from:

; L^2 is a displaceable group and R^{7a} is selected from the definition of R^7 or R^{22} above, and

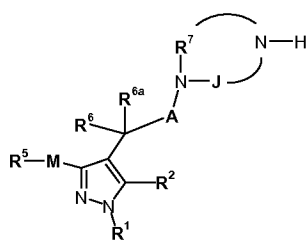
$L^2-R^{5''}$ is selected from:



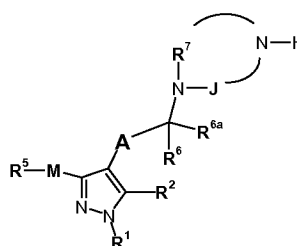
- (c) For compounds of Formula (I) wherein R^3 is a group of Formula (IIa), (IIb), (IIc) or (IId) and R^7 is other than part of a heterocyclic ring or hydrogen, reaction of a compound of Formula (I) wherein R^3 is a group of Formula (IIa), (IIb), (IIc) or (IId) and R^7 is hydrogen with a group of formula L^3-R^{7a} , wherein R^{7a} is as defined above for R^7 with the exclusion of hydrogen and L^3 is a displaceable group;
- (d) For compounds of Formula (I) wherein R^3 is a group of Formula (IIe) or (IIf) and R^{21} is other than hydrogen, reaction of a compound of Formula (I) wherein R^3 is a group of Formula (IIe) or (IIf) and R^{21} is hydrogen with a group of formula L^4-R^{21a} , wherein R^{21a} is as defined above for R^{21} with the exclusion of hydrogen and L^4 is a displaceable group;
- (e) For compounds of Formula (I) wherein R^3 is a group of Formula (IIe) or (IIf) and R^{22} is other than hydrogen, reaction of a compound of Formula (I) wherein R^3 is a group of Formula (IIe) or (IIf) and R^{22} is hydrogen with a group of formula L^5-R^{22a} , wherein R^{22a} is as defined above for R^{22} with the exclusion of hydrogen and L^5 is a displaceable group;
- (f) For compounds of Formula (I) wherein R^3 is a group of Formula (IIc) or (IId) and



the group together forms an optionally substituted nitrogen-containing heterocyclic ring containing 4-7 carbons atoms, reaction of a compound of Formula XXXIVa or XXXIVb, with a compound of Formula L^6-K-R^8 , wherein L^6 is a displaceable group

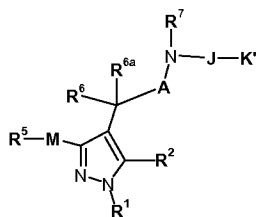


XXXIVa

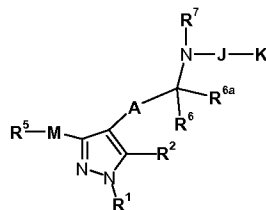


XXXIVb

- (g) For compounds of Formula (I) wherein R^3 is a group of Formula (IIc) or (IId), reaction of a compound of Formula XXXVa or XXXVb, with a compound of Formula $L^7-K''-R^8$, wherein L^7 is a displaceable group, and wherein the groups K' and K'' comprise groups which when reacted together form K ,



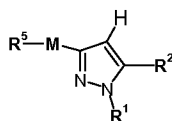
XXXVa



XXXVb

;

- (h) reaction of a compound of Formula XXXVI with a compound of the formula L^8-R^5 , wherein L^8 is a displaceable group

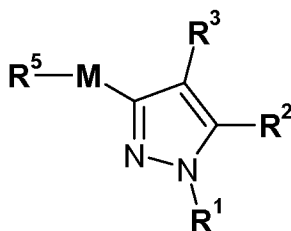


XXXVI ;

and thereafter if necessary:

- i) converting a compound of the Formula (I) into another compound of the Formula (I);
- ii) removing any protecting groups;
- iii) forming a salt, pro-drug or solvate.

17. (Currently Amended) A compound of Formula (I),



Formula (I),

or a salt, or in-vivo hydrolyzable ester pro-drug or solvate thereof, wherein

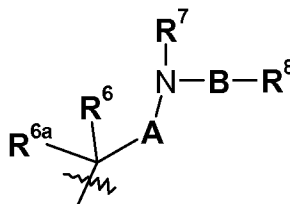
R¹ is hydrogen;

R² is 3,5-dimethylphenyl;

M is -CH₂-O-;

R⁵ is 2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy;

R³ is Formula (IIb),



Formula (IIb),

wherein,

R⁶ is hydrogen;

R^{6a} is methyl;

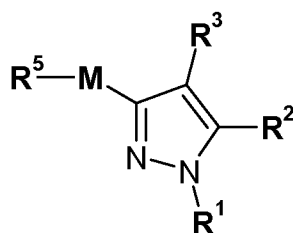
R⁷ is hydrogen;

R⁸ is 1,3-benzodioxol-5-yl;

A is methylene; and

B is selected from ethylene and butylene.

18. (New) A compound of Formula (I),



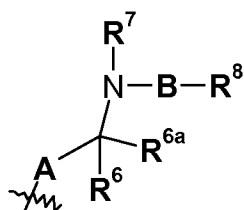
Formula (I)

wherein

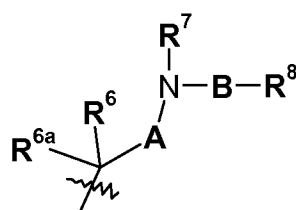
R¹ is selected from: hydrogen, optionally-substituted C₁₋₆alkyl, optionally substituted aryl or optionally-substituted arylC₁₋₆alkyl;

R² is optionally-substituted phenyl;

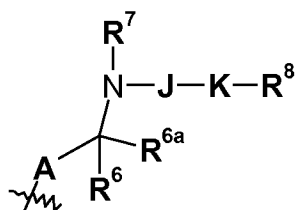
R³ is selected from a group of Formula (IIa) to Formula (IIf):



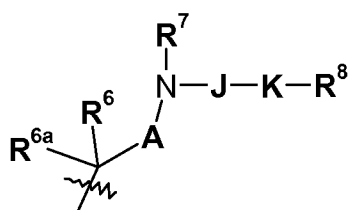
Formula (IIa)



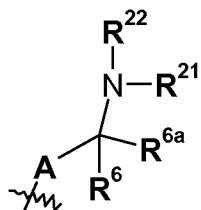
Formula (IIb)



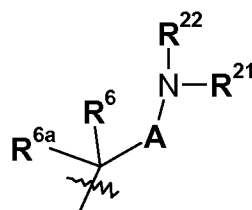
Formula (IIc)



Formula (IId)

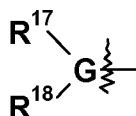


Formula (IIe)



Formula (IIf)

R⁵ is a group of Formula (III):



Formula (III)

R⁶ and **R^{6a}** are independently selected from hydrogen, fluoro, optionally substituted C₁₋₆alkyl or **R⁶** and **R^{6a}** taken together and the carbon atom to which they are attached form a carbocyclic ring of 3-7 atoms

R⁷ is selected from: hydrogen, optionally-substituted C₁₋₆alkyl, optionally-substituted arylC₁₋₆alkyl, optionally-substituted aryl, optionally substituted heterocyclyl, optionally substituted heterocyclylC₁₋₆alkyl, **R⁹**OC₁₋₆alkyl-, **R⁹R¹⁰**NC₁₋₆alkyl-, **R⁹R¹⁰**NC(O)C₁₋₆alkyl-, -C(NR⁹R¹⁰)=NH; or when **R³** is a group of Formula (IIc) or (IId) **R⁷** is of the formula -**J-K-R⁸**;

R⁸ is selected from:

- (i) hydrogen, C₁₋₆alkyl, C₂₋₆alkenyl, C₂₋₆alkynyl, haloC₁₋₆alkyl, C₁₋₄alkoxyC₁₋₄alkyl, hydroxy, hydroxyC₁₋₆alkyl, cyano, N-C₁₋₄alkylamino, N,N-di-C₁₋₄alkylamino, C₁₋₆alkyl-S(O)_n-, -O-**R^b**, -NR^bR^c, -C(O)-**R^b**, -C(O)O-**R^b**, -CONR^bR^c, NH-C(O)-**R^b** or -S(O)_nNR^bR^c, where **R^b** and **R^c** are independently selected from hydrogen and C₁₋₄alkyl optionally substituted with hydroxy, amino, N-C₁₋₄alkylamino, N,N-di-C₁₋₄alkylamino, HO-C₂₋₄alkyl-NH- or HO-C₂₋₄alkyl-N(C₁₋₄alkyl)-;
- (ii) nitro when **B** is a group of Formula (IV) and **X** is CH and **p** is 0;
- (iii) C₃₋₇cycloalkyl, aryl or arylC₁₋₆alkyl each of which is optionally substituted by **R¹²**, **R¹³** and **R¹⁴**;
- (iv) -(**Q**)-aryl, -(**Q**)-heterocyclyl, -aryl-(**Q**)-aryl, each of which is optionally substituted by **R¹²**, **R¹³** and **R¹⁴** wherein -(**Q**)- is selected from **E**, **F** or a direct bond;
- (v) heterocyclyl or heterocyclylC₁₋₆alkyl each of which is optionally substituted by up to 4 substituents independently selected from **R¹²**, **R¹³** and **R¹⁴**;
- (vi) a group selected from **R¹²**, **R¹³** and **R¹⁴**;

R⁹ and **R¹⁰** are independently selected from: hydrogen, hydroxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted arylC₁₋₆alkyl, an optionally substituted carbocyclic ring of 3-7 atoms,

optionally substituted heterocyclyl, optionally substituted heterocyclylC₁₋₆alkyl or **R**⁹ and **R**¹⁰ taken together can form an optionally substituted ring of 3-9 atoms or **R**⁹ and **R**¹⁰ taken together with the carbon atom to which they are attached form a carbonyl group;

R¹¹ is selected from: hydrogen, optionally substituted C₁₋₆alkyl, or N(**R**⁹**R**¹⁰);

R¹² is selected from: hydrogen, hydroxy, **R**¹⁷**R**¹⁸N(CH₂)_{cc}-, **R**¹⁷**R**¹⁸NC(O)(CH₂)_{cc}-, optionally substituted C₁₋₆alkyl- C(O)N(**R**⁹)(CH₂)_{cc}-, optionally substituted C₁₋₆alkyl-SO₂N(**R**⁹)-, optionally substituted aryl-SO₂N(**R**⁹)-, C₁₋₃perfluoroalkyl-SO₂N(**R**⁹)-, optionally substituted C₁₋₆alkyl-N(**R**⁹)SO₂-, optionally substituted aryl-N(**R**⁹)SO₂-, C₁₋₃perfluoroalkyl-N(**R**⁹)SO₂- optionally substituted C₁₋₆alkanoyl-N(**R**⁹)SO₂-, optionally substituted aryl-C(O)N(**R**⁹)SO₂-, optionally substituted C₁₋₆alkyl-S(O_n) -, optionally substituted aryl-S(O_n) -, C₁₋₃perfluoroalkyl-, C₁₋₃perfluoroalkoxy, optionally substituted C₁₋₆alkoxy, carboxy, halo, nitro or cyano;

R¹³ and **R**¹⁴ are independently selected from: hydrogen, hydroxy, oxo, optionally substituted C₁₋₆alkyl, optionally substituted C₁₋₆alkanoyl, optionally substituted C₂₋₆alkenyl, cyano, nitro, C₁₋₃perfluoroalkyl-, C₁₋₃perfluoroalkoxy, optionally substituted aryl, optionally substituted arylC₁₋₆alkyl, **R**⁹O(CH₂)_s-, **R**⁹(O)O(CH₂)_s-, **R**⁹OC(O)(CH₂)_s-, **R**¹⁶S(O_n)(CH₂)_s-, **R**⁹**R**¹⁰NC(O)(CH₂)_s- or halo;

R¹⁵ is selected from: hydrogen, optionally substituted C₁₋₆alkyl, **R**¹⁹OC(O)-, **R**⁹**R**¹⁰NC(O)-, **R**⁹C(O)-, **R**⁹S(O_n)-;

R¹⁶ is selected from: hydrogen, C₁₋₆alkyl, C₁₋₃perfluoroalkyl or optionally-substituted aryl;

R¹⁷ is independently selected from: hydrogen, hydroxy, cyano or optionally substituted C₁₋₆alkyl;

R¹⁸ is a group of formula **R**^{18a}-C(**R**⁹**R**¹⁰)₀₋₁- wherein **R**^{18a} is selected from: **R**¹⁹OC(O)-, **R**⁹**R**¹⁰NC(O)-, **R**⁹**R**¹⁰N-, **R**⁹C(O)-, **R**⁹C(O)N(**R**¹⁰)-, **R**⁹**R**¹⁰NC(O)-, **R**⁹**R**¹⁰NC(O)N(**R**¹⁰)-, **R**⁹SO₂N(**R**¹⁰)-, **R**⁹**R**¹⁰NSO₂N(**R**¹⁰)-,

$\mathbf{R}^9\text{C}(\text{O})\text{O}-$, $\mathbf{R}^9\text{OC}(\text{O})-$, $\mathbf{R}^9\mathbf{R}^{10}\text{NC}(\text{O})\text{O}-$, $\mathbf{R}^9\text{O}-$, $\mathbf{R}^9\text{S}(\text{O}_n)-$, $\mathbf{R}^9\mathbf{R}^{10}\text{NS}(\text{O}_n)-$,

hydrogen, optionally substituted C_{1-6} alkyl, optionally substituted

heterocyclyl;

or \mathbf{R}^{17} and \mathbf{R}^{18} when taken together form an optionally substituted carbocyclic ring of 3-7 atoms or optionally substituted heterocyclyl;

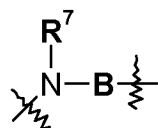
\mathbf{R}^{19} is selected from: hydrogen, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted C_{3-7} cycloalkyl, optionally substituted heterocyclyl or optionally substituted heterocyclyl C_{1-6} alkyl;

\mathbf{R}^{21} and \mathbf{R}^{22} are independently selected from hydrogen, optionally substituted C_{1-6} alkyl, optionally substituted C_{3-7} cycloalkyl, optionally substituted heterocyclyl, optionally substituted heterocyclyl C_{1-6} alkyl, optionally substituted C_{3-6} alkenyl, optionally substituted C_{3-6} alkynyl, -
 $(\text{C}_{1-5}\text{alkyl})_{\text{aa}}-\text{S}(\text{O}_n)-(\text{C}_{1-5}\text{alkyl})_{\text{bb}}-$; $\mathbf{R}^9\mathbf{R}^{10}\text{NC}_{2-6}\text{alkyl}$, $\mathbf{R}^9\text{OC}_{2-6}\text{alkyl}$ or $\mathbf{R}^9\mathbf{R}^{10}\text{NC}(\text{O})\text{C}_{2-6}\text{alkyl}$, with the proviso that \mathbf{R}^9 and \mathbf{R}^{10} independently or taken together are not optionally substituted aryl or optionally substituted aryl C_{1-6} alkyl; or

\mathbf{R}^{21} and \mathbf{R}^{22} taken together form an optionally substituted non-aromatic heterocyclic ring;

A is selected from a direct bond, optionally substituted C_{1-5} alkylene, carbonyl or $-\text{C}(\text{O})-\text{C}(\mathbf{R}^d\mathbf{R}^d)-$, wherein \mathbf{R}^d is independently selected from a direct bond hydrogen and C_{1-2} alkyl;

B is C_{1-6} alkylene, C_{3-6} alkenylene, $-(\text{C}_{1-5}\text{alkyl})_{\text{aa}}-\text{O}-(\text{C}_{1-5}\text{alkyl})_{\text{bb}}-$, $-(\text{C}_{1-5}\text{alkyl})_{\text{aa}}-\text{C}(\text{O})-(\text{C}_{1-5}\text{alkyl})_{\text{bb}}-$, $-(\text{CH}_2)_{\text{s1}}-\text{C}(\text{O})\text{N}(\mathbf{R}^9)-$, or the group



forms an optionally substituted saturated C_{4-7} heterocyclic ring,

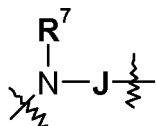
wherein **aa** and **bb** are independently 0 or 1 and wherein the combined length of $(\text{C}_{1-5}\text{alkyl})_{\text{aa}}$, $(\text{C}_{1-5}\text{alkyl})_{\text{bb}}$ is less than or equal to C_5 alkyl and wherein C_{1-6} alkylene is optionally substituted by hydroxy.

E is $-O-$, $-S(O_n)-$, $-C(O)-$, $-NR^{15}-$ or $-C(R^9R^{10})_q-$;

F is $-E(CH_2)_r-$;

G is selected from: hydrogen, halo, N, O, $S(O_n)$, $C(O)$, $C(R^9R^{10})_t$, optionally substituted C_{2-6} alkenylene, optionally substituted C_{2-6} alkynylene or a direct bond to R^{18} ,

J is a group of the formula: $-(CH_2)_s-L-(CH_2)_s-$ wherein when **s** is greater than 0, the alkylene group is optionally substituted,



or the group together forms an optionally substituted heterocyclic ring containing 4-7 carbons atoms;

K is selected from: a direct bond, $-(CH_2)_{s1}-$, $-(CH_2)_{s1}-O-(CH_2)_{s2}-$, $-(CH_2)_{s1}C(O)-(CH_2)_{s2}-$, $-(CH_2)_{s1}S(O_n)-(CH_2)_{s2}-$, $-(CH_2)_{s1}N(R^{18})-(CH_2)_{s2}-$, $-(CH_2)_{s1}-C(O)N(R^9)-(CH_2)_{s2}-$, $-(CH_2)_{s1}-N(R^9)C(O)-(CH_2)_{s2}-$, $-(CH_2)_{s1}-N(R^9)C(O)N(R^9)-(CH_2)_{s2}-$, $-(CH_2)_{s1}-OC(O)-(CH_2)_{s2}-$, $-(CH_2)_{s1}-C(O)O-(CH_2)_{s2}-$, $-(CH_2)_{s1}-N(R^9)C(O)O-(CH_2)_{s2}-$, $-(CH_2)_{s1}-OC(O)N(R^9)-(CH_2)_{s2}-$, $-(CH_2)_{s1}-OS(O_n)-(CH_2)_{s2}-$, or $-(CH_2)_{s1}-S(O_n)-O-(CH_2)_{s2}-$, $-(CH_2)_{s1}-S(O)_2N(R^9)-(CH_2)_{s2}-$, $-(CH_2)_{s1}-N(R^9)S(O)_2-(CH_2)_{s2}-$; wherein the $-(CH_2)_{s1}-$ and $-(CH_2)_{s2}-$ groups are independently optionally substituted by hydroxy or C_{1-4} alkyl;

L is selected from optionally substituted aryl or optionally substituted heterocyclyl;

M is $-(CH_2)-O-$;

n is an integer from 0 to 2;

p is an integer from 0 to 4;

q is an integer from 0 to 4;

r is an integer from 0 to 4;

s is an integer from 0 to 4;

s1 and **s2** are independently selected from an integer from 0 to 4, and

s1+s2 is less than or equal to 4;

t is an integer from 0 to 4;

aa and **bb** are independently 0 or 1; and

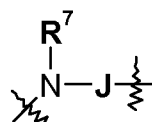
cc is an integer between 0 to 2;

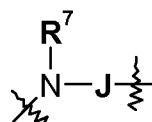
with the proviso that

- (i) when **G** is hydrogen or halo, then **R**¹⁷ and **R**¹⁸ are both absent;
- (ii) when **G** is O, S(O_n), C(O) or C(**R**^{11**R**¹²)_t then **G** is substituted by a single group independently selected from the definition of **R**¹⁷ or **R**¹⁸ and when **G** is a direct bond to **R**¹⁸ then **G** is substituted by a single group selected from **R**¹⁸;}
- (iii) when **R**³ is a group of Formula (IIb), **B** is a group of Formula (IV), **R**⁸ is selected from group (i) or (ii) above, **R**¹¹ is a group of the formula N(**R**^{10**R**¹¹) and **R**¹, **R**² and **R**⁵ are as defined above then **R**⁴ cannot be hydrogen;}
- (iv) **R**³ cannot be unsubstituted pyridyl or unsubstituted pyrimidinyl; and
- (v) when **R**³ is pyrazolyl substituted by phenyl or pyrazolyl substituted by phenyl and acetyl, **R**⁵-**M** is hydroxyl or acetyloxy, **R**² is unsubstituted phenyl, then **R**¹ cannot be hydrogen or acetyl;

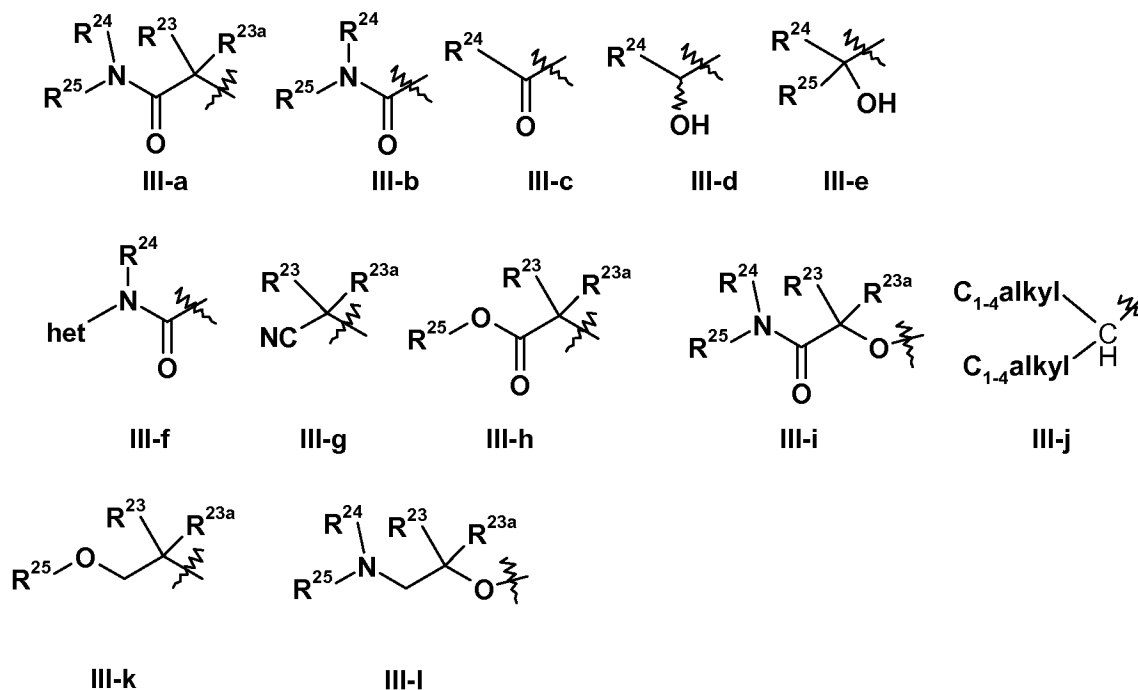
or a salt, or in-vivo hydrolyzable ester thereof.

- 19. (New) The compound of Claim 18, wherein **R**¹ is hydrogen.
- 20. (New) The compound of Claim 18, wherein **R**³ is selected from a group of Formula (IIa) or Formula (IIb).
- 21. (New) The compound of Claim 20, wherein **B** is optionally substituted C₁₋₆alkylene.
- 22. (New) The compound of Claim 18, wherein **R**³ is selected from a group of Formula (IIc) or Formula (IId).



23. (New) The compound of Claim 22 wherein the group  together forms an optionally substituted heterocyclic ring containing 4-7 carbons atoms.
24. (New) The compound according to Claim 23 wherein **K** is selected from:
 -(CH₂)_s-, -(CH₂)_s-O-(CH₂)_s-, -(CH₂)_s-C(O)-(CH₂)_s-, -(CH₂)_s-N(**R**¹⁸)-(CH₂)_s-,
 -(CH₂)_s-C(O)N(**R**¹⁸)-(CH₂)_s-, -(CH₂)_s-N(**R**¹⁸)C(O)-(CH₂)_s-,
 -(CH₂)_s-S(O)₂N(**R**¹⁸)-(CH₂)_s-, or -(CH₂)_s-NHS(O)₂-(CH₂)_s-.
25. (New) The compound of Claim 20 wherein **R**⁸ is selected from:
- (i) hydrogen, C₁₋₆alkyl, C₂₋₆alkenyl, haloC₁₋₆alkyl, hydroxy, cyano, C₁₋₆alkylS(O)_n-, -O-**R**^b, C₁₋₄alkoxyC₁₋₄alkyl, -C(O)-**R**^b, C(O)O-**R**^b, -NH-C(O)-**R**^b, N,N-di-C₁₋₄alkylamino, -S(O)_nNR^b**R**^c where **R**^b and **R**^c are independently selected from hydrogen and C₁₋₆alkyl, and **n** is 0, 1 or 2;
 - (ii) -(**Q**)-aryl, optionally substituted by up to 3 groups selected from **R**¹², **R**¹³ and **R**¹⁴;
 - (iii) C₄₋₇heterocyclyl, optionally substituted by up to 3 groups selected from **R**¹², **R**¹³ and **R**¹⁴, or
 - (iv) C₃₋₇carbocyclyl, optionally substituted by up to 3 groups selected from **R**¹², **R**¹³ and **R**¹⁴.

26. (New) The compound of Claim 18 wherein R^5 is a group of Formula (III) wherein the group of Formula (III) is selected from any one of **III-a** to **III-l**;



wherein:

het represents an optionally substituted 3- to 8- membered heterocyclic ring containing from 1 to 4 heteroatoms independently selected from O, N and S;
 R^{23} and R^{23a} are independently selected from hydrogen, fluoro or optionally substituted C_{1-8} alkyl; or R^{23} and R^{23a} together with the carbon to which they are attached form an optionally substituted 3 to 7-membered cycloalkyl ring
 R^{24} is selected from hydrogen, optionally substituted C_{1-8} alkyl, optionally substituted aryl, $-R^d$ -Ar, where R^d represents C_{1-8} alkylene and Ar represents optionally substituted aryl, and optionally substituted 3- to 8-membered heterocyclic ring optionally containing from 1 to 3 further heteroatoms independently selected from O, N and S;
 R^{25} is selected from hydrogen; optionally substituted C_{1-8} alkyl and optionally substituted aryl;

or where the group of Formula (III) represents a group of Formula **III-a**, **III-b** or **III-i**, then the group $\text{NR}^{24}(-\text{R}^{25})$ represents an optionally substituted 3- to 8- membered heterocyclic ring optionally containing from 1 to 3 further heteroatoms independently selected from O, N and S;

or where the group of Formula (III) represents structure **III-e**, R^{24} and R^{25} together with the carbon to which they are attached represents an optionally substituted 3- to 8- membered heterocyclic ring optionally containing from 1 to 4 heteroatoms independently selected from O, N and S.

27. (New) The compound of Claim 18 wherein the optional substituents on R^2 are selected from cyano, NR^eR^f , optionally substituted C_{1-8} alkyl, optionally substituted C_{1-8} alkoxy or halo wherein R^e and R^f are independently selected from hydrogen, C_{1-6} alkyl or aryl.

28. (New) The compound of Claim 18 selected from:

2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-pyrid-4-ylethyl]-(2*S*)-propylamine;
2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-pyrid-4-ylbutyl]-(2*S*)-propylamine;
2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[4-(4-methoxyphenyl)butyl]-(2*S*)-propylamine;
2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-(43-trifluoromethylphenyl)ethyl]-(2*S*)-propylamine;
2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-(4-fluorophenyl)ethyl]-(2*S*)-propylamine;

2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-(3-methoxyphenyl)ethyl]-(2*S*)-propylamine;
2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-(4-methoxyphenyl)ethyl]-(2*S*)-propylamine;
2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.1]heptan-7-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-(4-methylsulphonylaminophenyl)ethyl]-(2*S*)-propylamine; and
2-[3-(2,2-dimethyl-3-oxo-3-{azabicyclo[2.2.2]oct-2-yl}propoxy)-5-(3,5-dimethylphenyl)-1*H*-pyrazol-4-yl]-*N*-[2-(1,3-benzodioxol-5-yl)ethyl]-(2*S*)-propylamine;
or a salt, or in-vivo hydrolyzable ester thereof.

29. (New) A pharmaceutical formulation comprising a compound, or salt, or in-vivo hydrolyzable ester thereof, according to Claim 18 and a pharmaceutically acceptable diluent or carrier.